

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
 Data File : BG055792.D
 Acq On : 25 Nov 2022 22:48
 Operator : CG/JU
 Sample : N5641-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SOIL-PILE-10

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055792.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.286	334	340	349	rBV	91945	146859	7.37%	1.026%
2	5.768	414	422	439	rBV	601147	988380	49.60%	6.903%
3	7.384	689	697	713	rBV	596578	1061740	53.28%	7.416%
4	8.236	834	842	849	rBV	162815	283041	14.20%	1.977%
5	9.405	1033	1041	1051	rBV	375172	655644	32.90%	4.579%
6	10.803	1272	1279	1299	rBV	77472	179983	9.03%	1.257%
7	11.062	1315	1323	1330	rBV	219934	400649	20.11%	2.798%
8	13.482	1727	1735	1743	rBV	965761	1493847	74.96%	10.434%
9	14.857	1961	1969	1975	rBV	357599	545001	27.35%	3.806%
10	16.338	2214	2221	2234	rBV	707045	1126570	56.53%	7.868%
11	17.325	2384	2389	2394	rBV2	73997	94912	4.76%	0.663%
12	17.431	2402	2407	2411	rBV2	25518	33998	1.71%	0.237%
13	17.601	2429	2436	2441	rBV	414790	642812	32.26%	4.490%
14	17.642	2441	2443	2449	rVB	27569	32689	1.64%	0.228%
15	18.230	2539	2543	2549	rVB	19140	25135	1.26%	0.176%
16	18.453	2577	2581	2590	rBV3	71258	113720	5.71%	0.794%
17	19.634	2778	2782	2787	rVB	71228	104581	5.25%	0.730%
18	19.716	2792	2796	2803	rBV3	28402	44873	2.25%	0.313%
19	19.998	2840	2844	2851	rVB	76070	112777	5.66%	0.788%
20	20.186	2870	2876	2882	rBV	1517424	1992752	100.00%	13.918%
21	20.468	2921	2924	2932	rVB5	16073	33846	1.70%	0.236%
22	20.750	2969	2972	2975	rBV	24196	25588	1.28%	0.179%
23	20.780	2975	2977	2982	rVB3	17704	21943	1.10%	0.153%
24	20.862	2987	2991	2997	rVB	248757	315043	15.81%	2.200%
25	21.297	3062	3065	3070	rVB	59752	78307	3.93%	0.547%
26	21.502	3095	3100	3104	rBV7	29412	76213	3.82%	0.532%
27	21.890	3159	3166	3172	rBV2	403795	746085	37.44%	5.211%
28	22.695	3298	3303	3308	rBV3	80349	156729	7.86%	1.095%
29	22.760	3309	3314	3317	rBV2	146125	291096	14.61%	2.033%
30	22.912	3334	3340	3343	rBV2	269846	593782	29.80%	4.147%
31	22.960	3344	3348	3353	rVV	329777	676810	33.96%	4.727%
32	23.212	3387	3391	3397	rBV	246116	459904	23.08%	3.212%
33	25.298	3737	3746	3757	rVB2	251575	762467	38.26%	5.325%

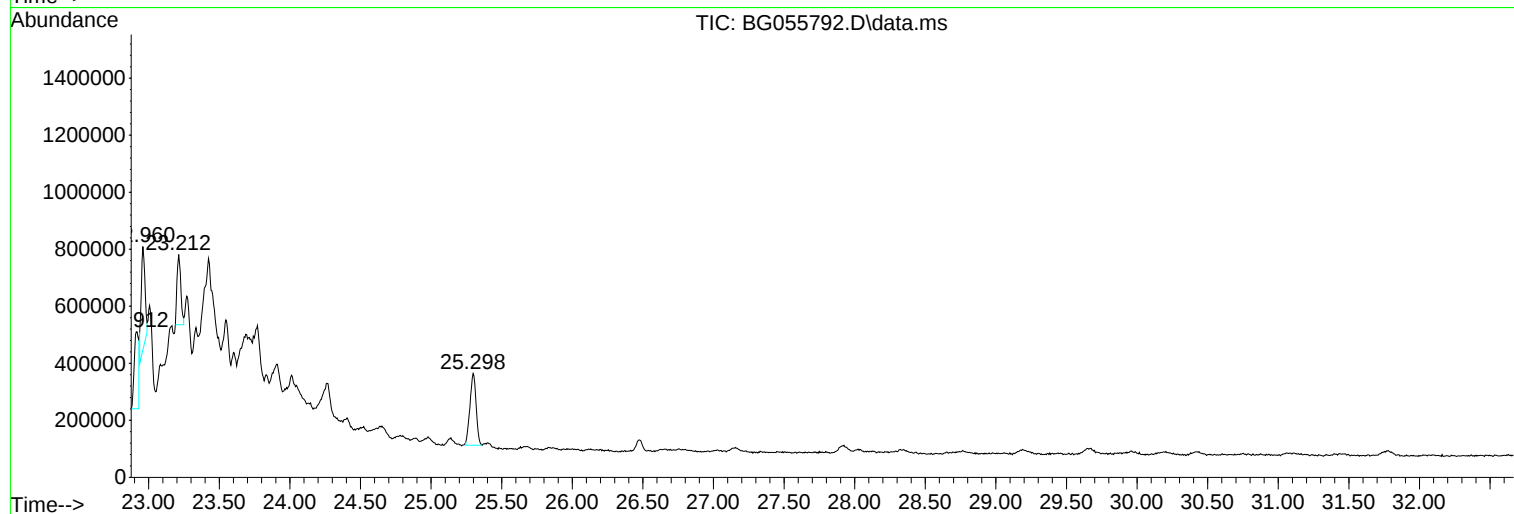
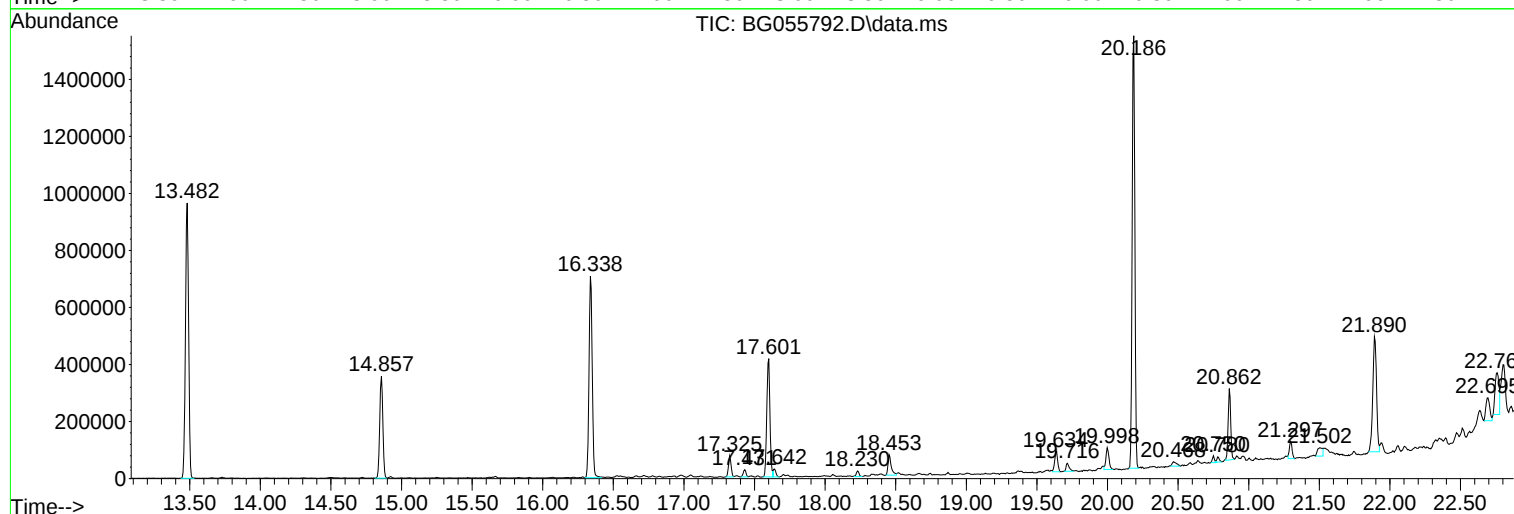
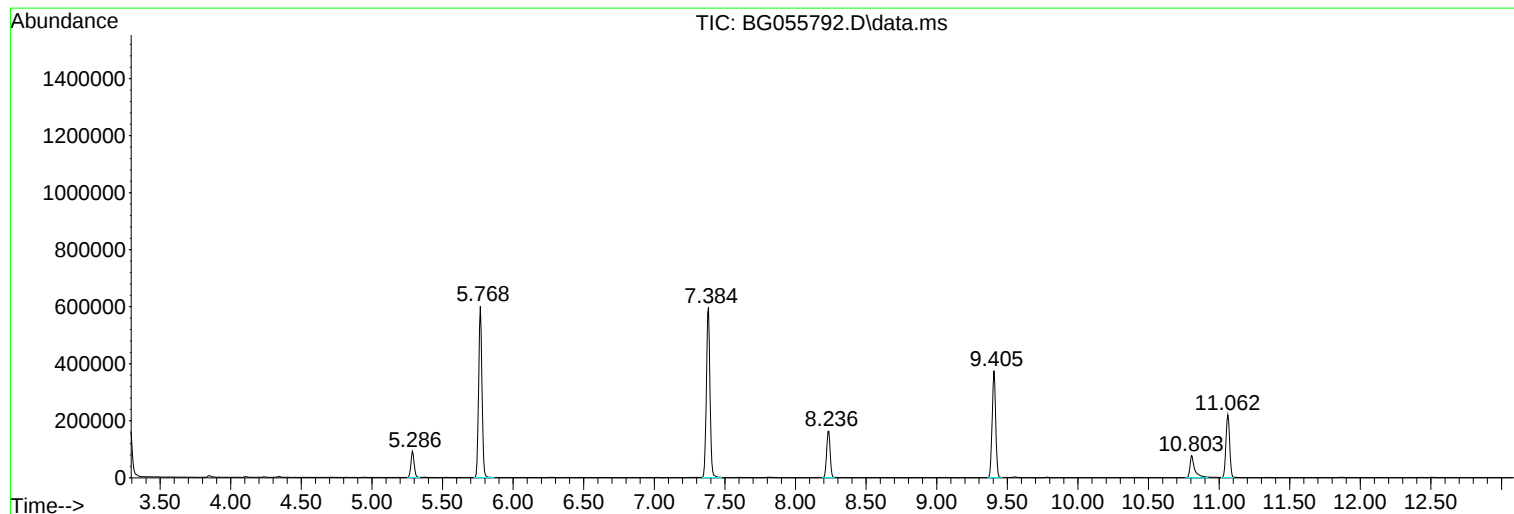
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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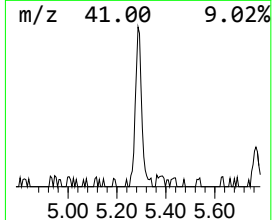
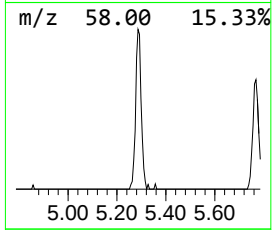
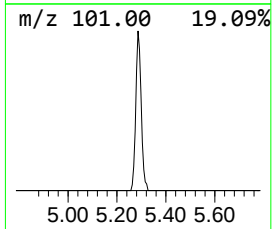
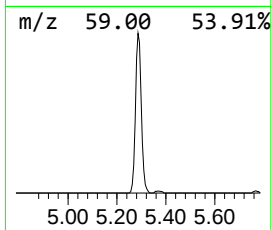
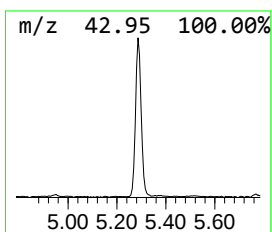
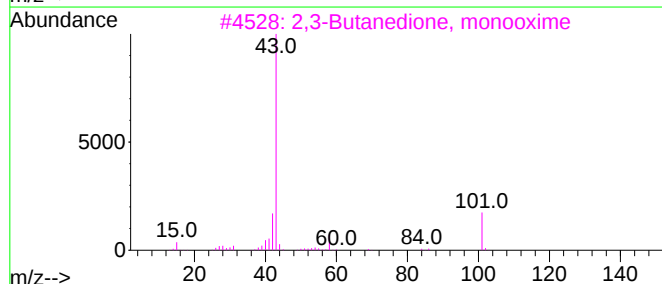
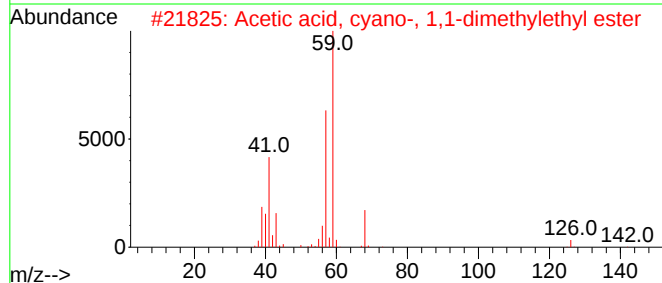
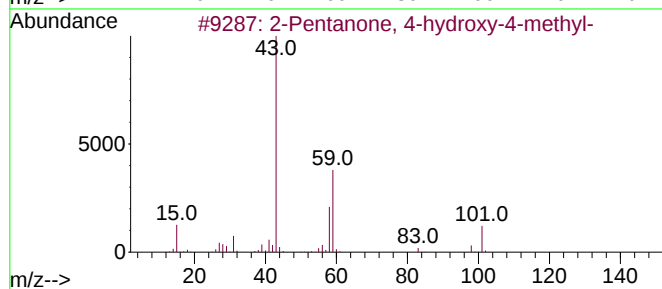
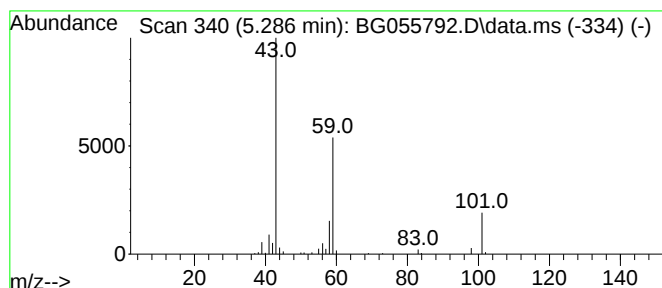
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.286	10.38 ng	146859	1,4-Dichlorobenzene-d4	8.236

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			3-Heptanol	116	C7H16O	000589-82-2	9



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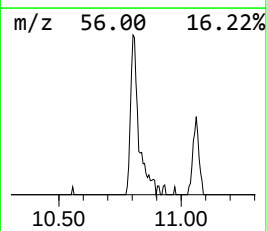
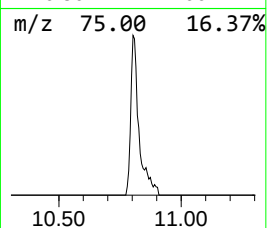
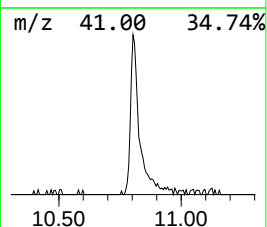
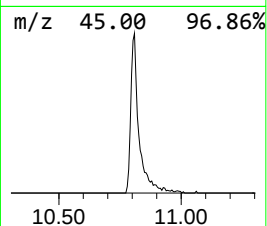
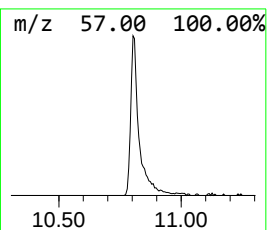
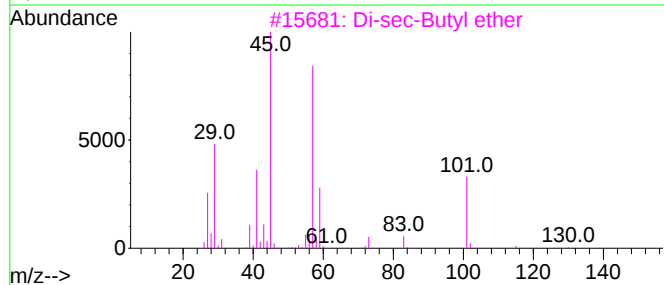
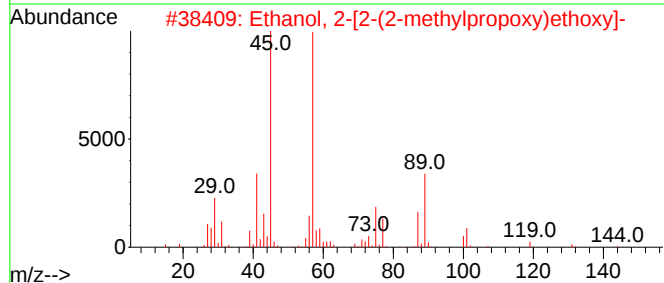
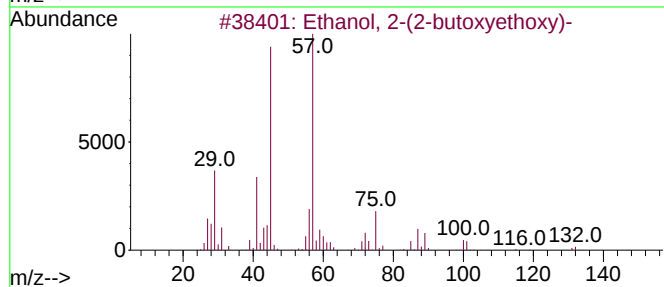
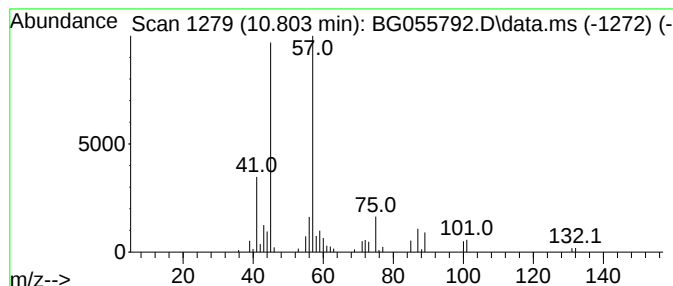
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG11622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.803	8.98 ng	179983	Naphthalene-d8	11.062	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78
3	Di-sec-Butyl ether	130	C8H18O	006863-58-7	50
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	47
5	3,6,9,12-Tetraoxatetradeca-1,13-...	202	C10H18O4	000765-12-8	45



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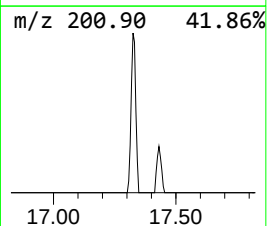
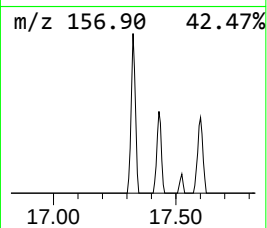
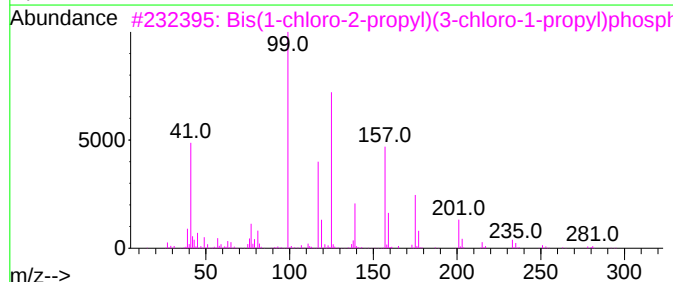
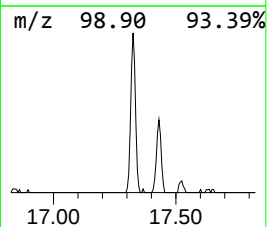
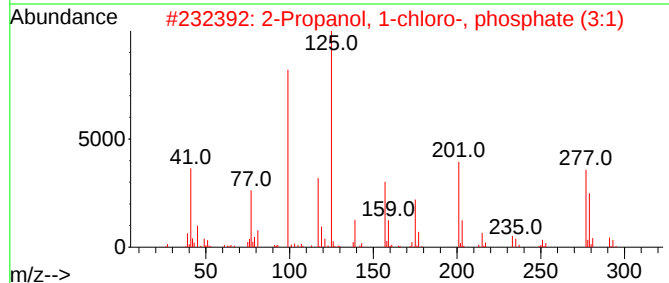
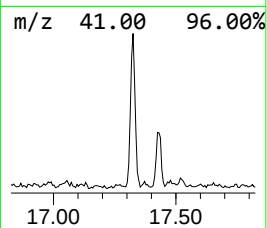
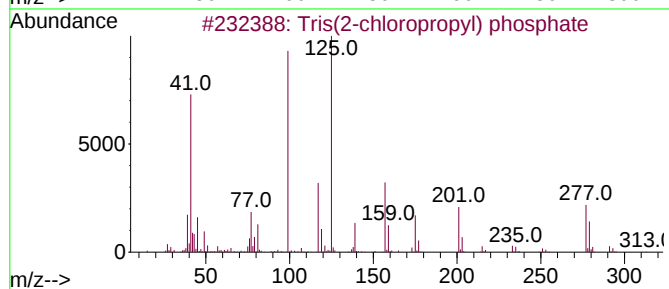
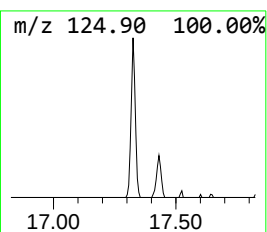
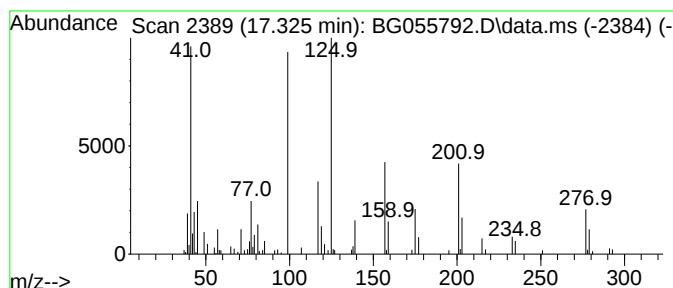
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TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Tris(2-chloropropyl) phosphate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.325	2.95 ng	94912	Phenanthrene-d10		17.601	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tris(2-chloropropyl) phosphate		326	C9H18Cl3O4P	006145-73-9	90
2	2-Propanol, 1-chloro-, phosphate...		326	C9H18Cl3O4P	013674-84-5	90
3	Bis(1-chloro-2-propyl)(3-chloro-...		326	C9H18Cl3O4P	137909-40-1	52
4	Bis(3-chloro-1-propyl)(1-chloro-...		326	C9H18Cl3O4P	137888-35-8	38
5	Tris(3-chloropropyl) phosphate		326	C9H18Cl3O4P	001067-98-7	35



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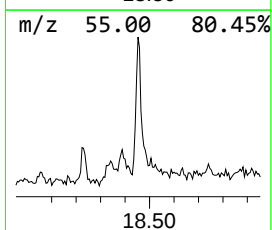
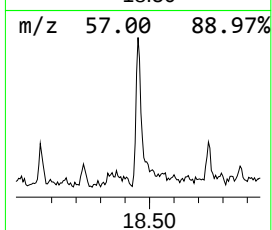
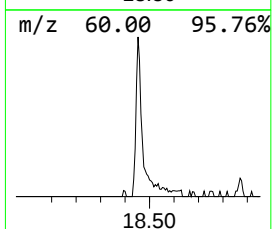
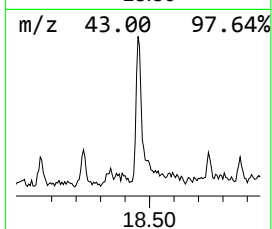
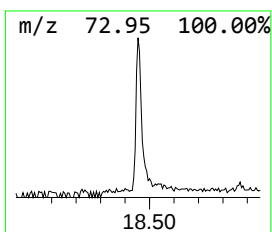
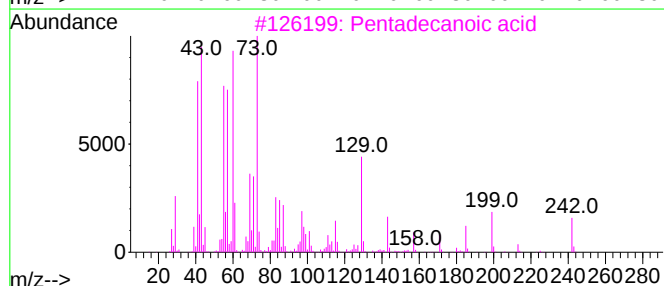
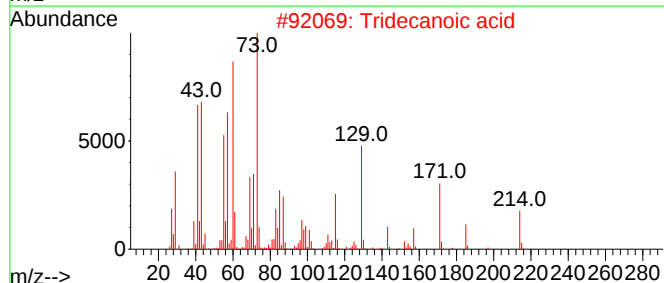
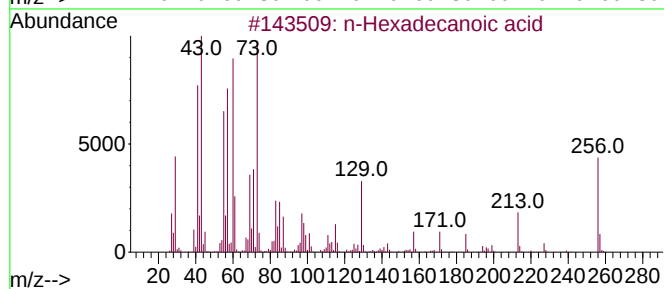
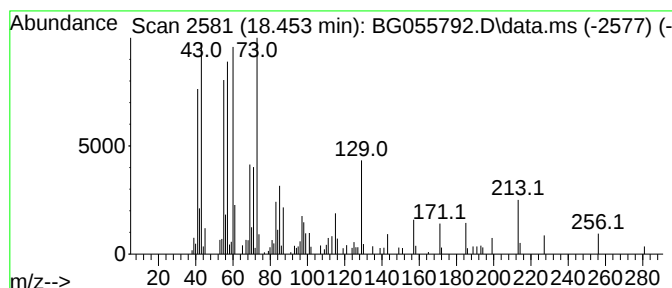
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.453	3.54 ng	113720	Phenanthrene-d10	17.601	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5	n-Decanoic acid	172	C10H20O2	000334-48-5	81



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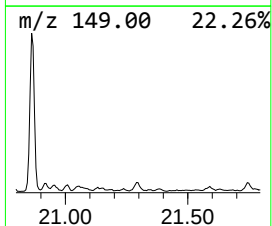
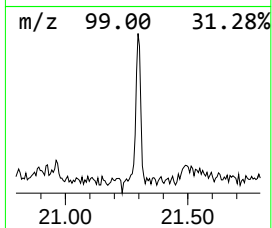
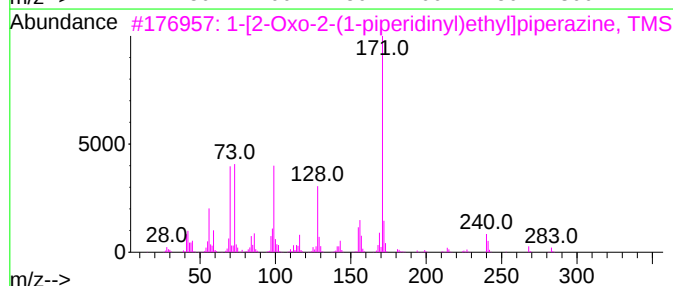
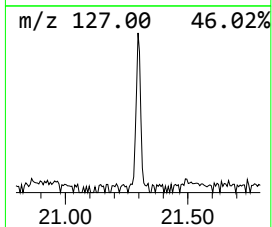
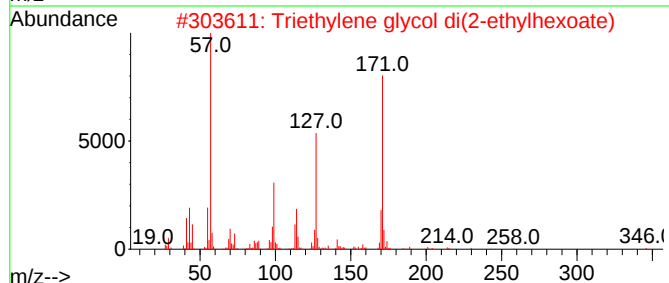
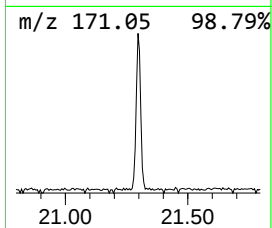
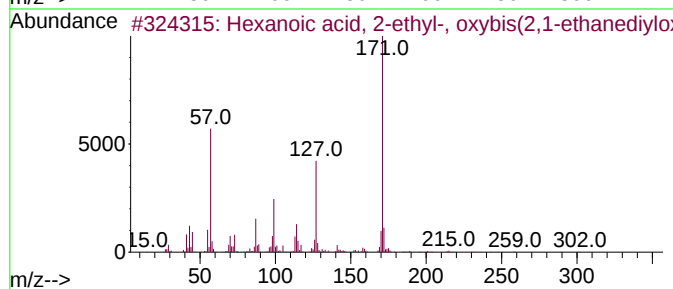
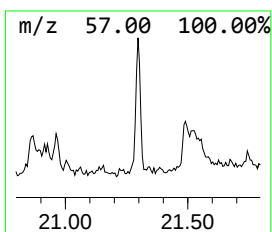
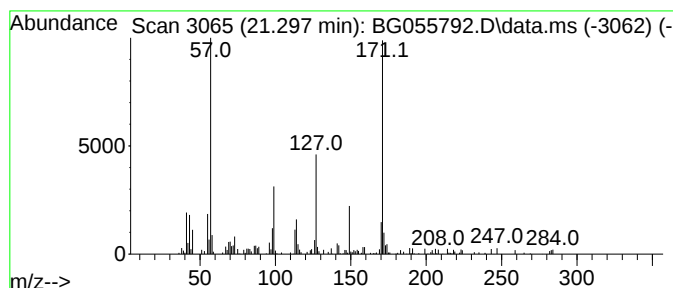
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Hexanoic acid, 2-ethyl-, ox... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.297	2.10 ng	78307	Chrysene-d12	21.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-, oxybis(...	446	C24H46O7	018268-70-7	80
2			Triethylene glycol di(2-ethylhex...	402	C22H42O6	000094-28-0	64
3			1-[2-Oxo-2-(1-piperidinyl)ethyl]...	283	C14H29N3OSi	1000474-15-5	43
4			3H,6H-Thieno[3,4-c]isoxazole, 3a...	171	C8H13NOS	127865-49-0	35
5			2,6-Difluorophenyl isocyanate	171	C7H3F2NS	065295-69-4	35



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Sample : N5641-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SOIL-PILE-10

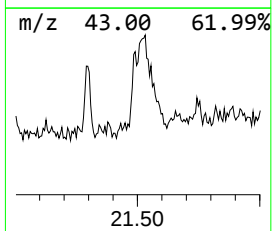
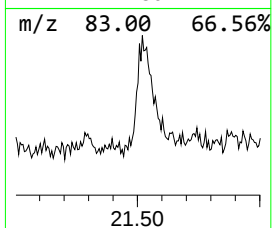
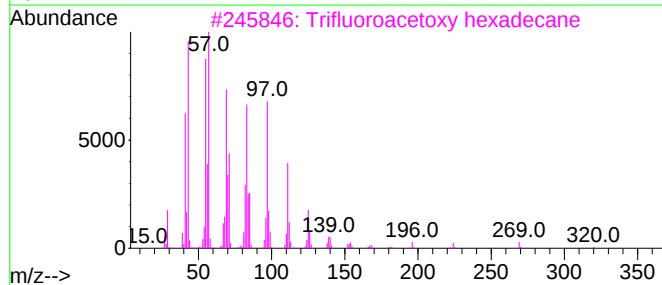
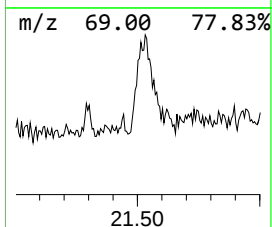
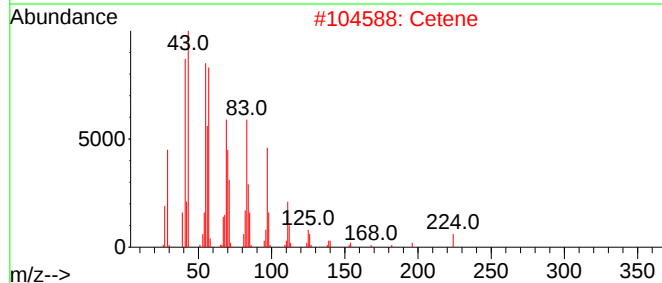
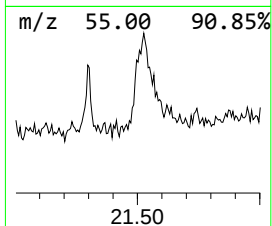
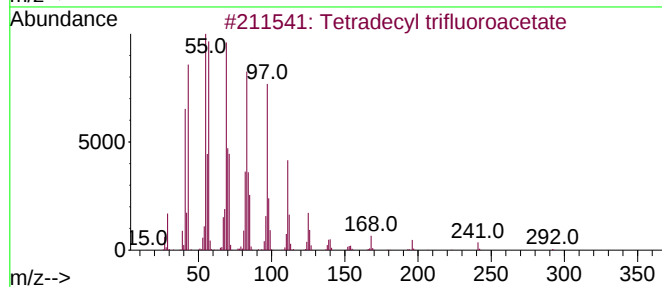
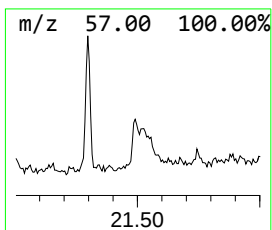
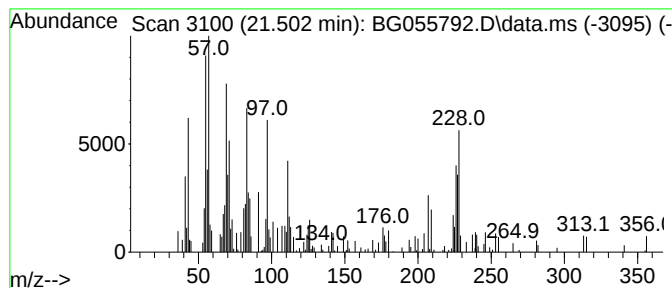
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown21.502 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.502	2.04 ng	76213	Chrysene-d12	21.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecyl trifluoroacetate	310	C16H29F3O2	006222-02-2	45
2			Cetene	224	C16H32	000629-73-2	44
3			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	43
4			1-Tetracosene	336	C24H48	010192-32-2	43
5			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
Data File : BG055792.D
Acq On : 25 Nov 2022 22:48
Operator : CG/JU
Sample : N5641-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SOIL-PILE-10

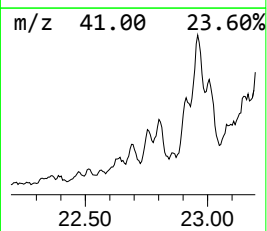
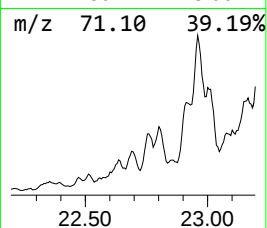
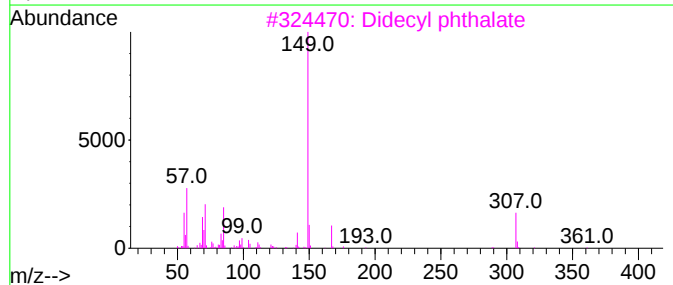
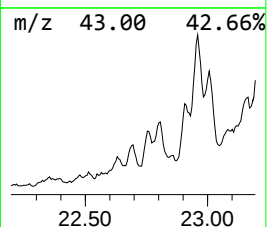
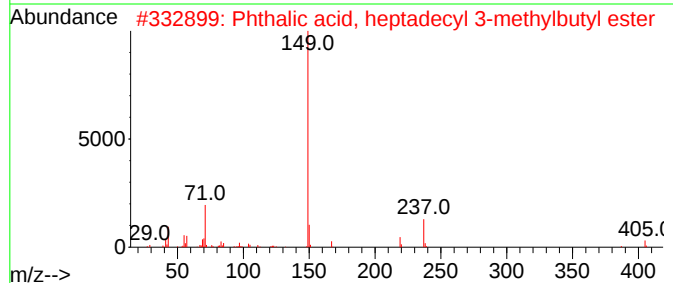
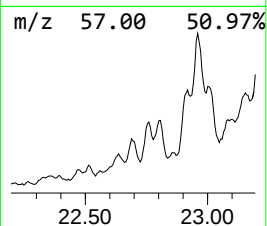
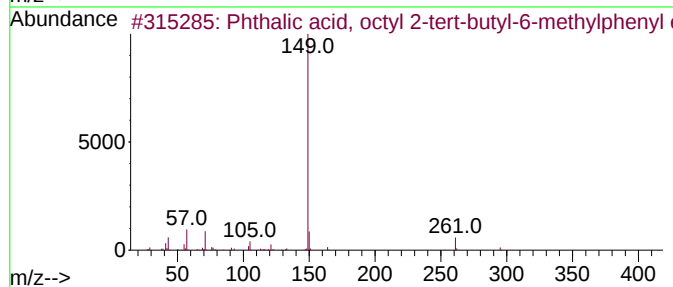
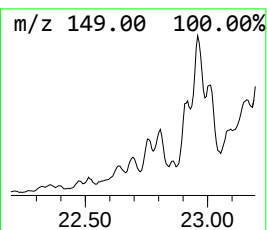
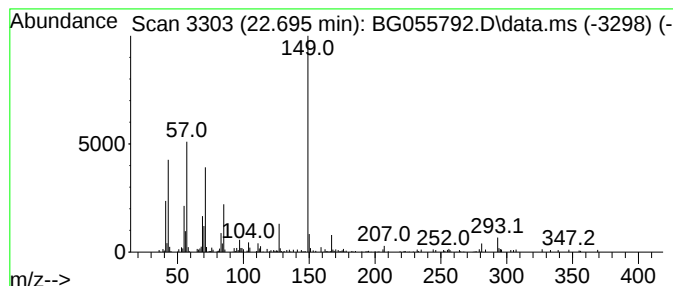
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Phthalic acid, octyl 2-tert... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.695	4.20 ng	156729	Chrysene-d12	21.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, octyl 2-tert-buty...	424	C27H36O4	1000357-09-8	53
2			Phthalic acid, heptadecyl 3-meth...	474	C30H50O4	1000356-81-8	53
3			Didecyl phthalate	446	C28H46O4	000084-77-5	53
4			Phthalic acid, butyl 2-methylpen...	306	C18H26O4	1000356-90-8	43
5			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
Data File : BG055792.D
Acq On : 25 Nov 2022 22:48
Operator : CG/JU
Sample : N5641-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

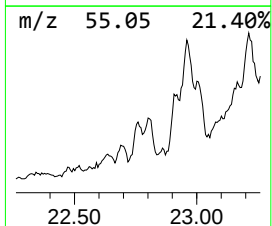
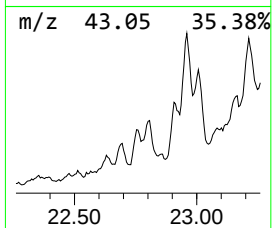
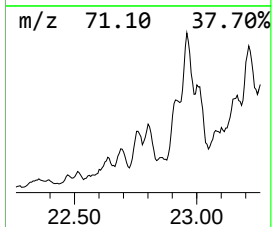
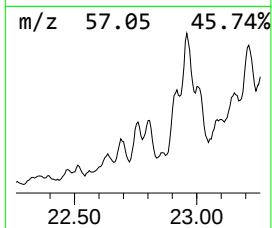
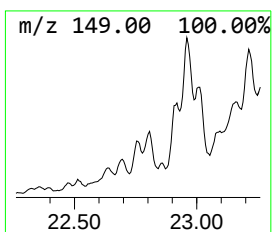
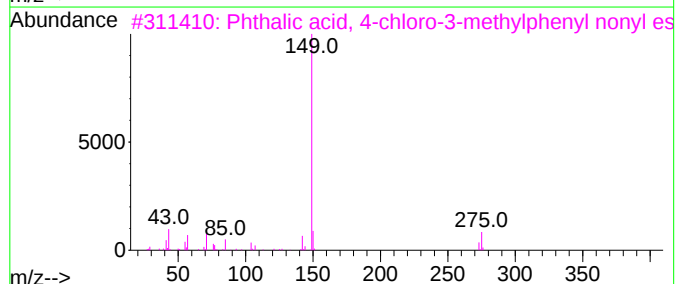
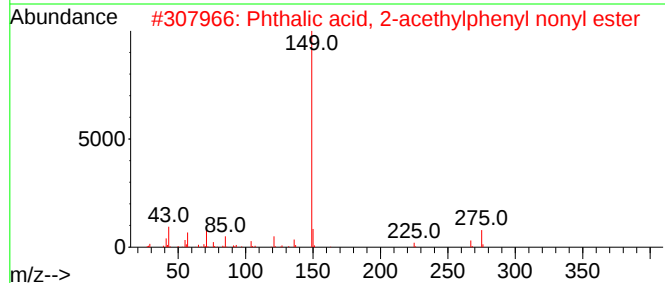
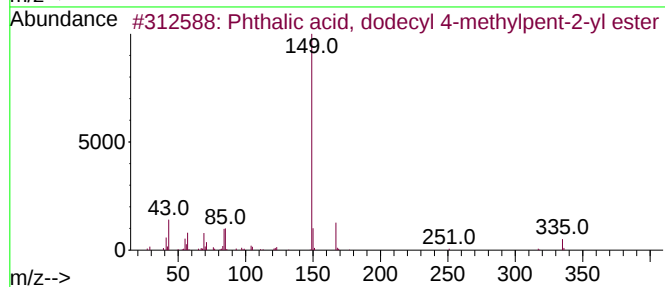
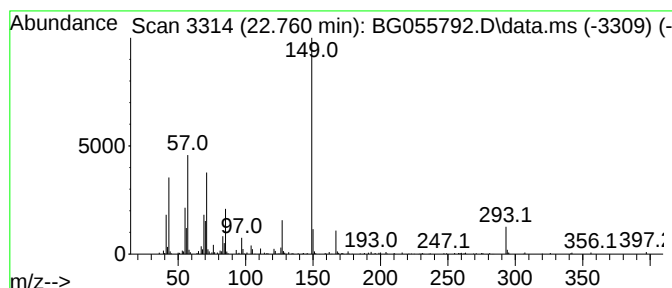
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG11622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phthalic acid, dodecyl 4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.760	7.80 ng	291096	Chrysene-d12	21.890	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, dodecyl 4-methylp...	418	C26H42O4	1000356-88-2	59
2	Phthalic acid, 2-acethylphenyl n...	410	C25H30O5	1000315-82-0	59
3	Phthalic acid, 4-chloro-3-methyl...	416	C24H29ClO4	1000309-85-0	59
4	Phthalic acid, 3,4-dichloropheny...	394	C20H20Cl2O4	1000315-27-5	53
5	Phthalic acid, isohexyl 3-methyl...	320	C19H28O4	1000356-80-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
Data File : BG055792.D
Acq On : 25 Nov 2022 22:48
Operator : CG/JU
Sample : N5641-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SOIL-PILE-10

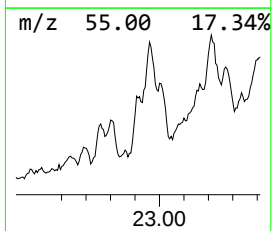
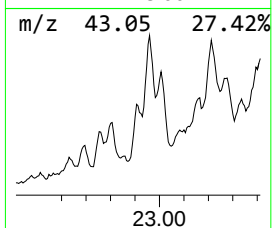
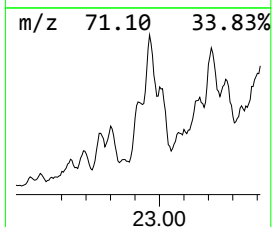
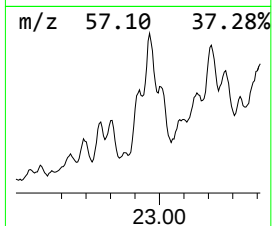
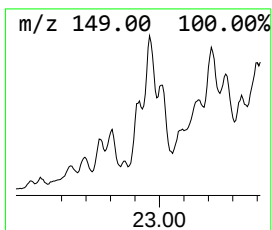
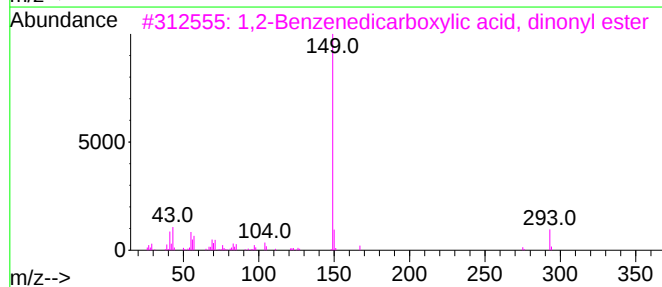
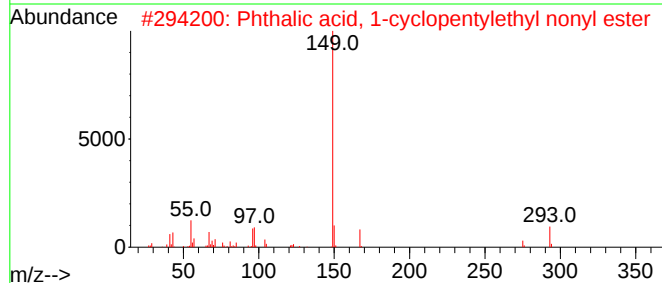
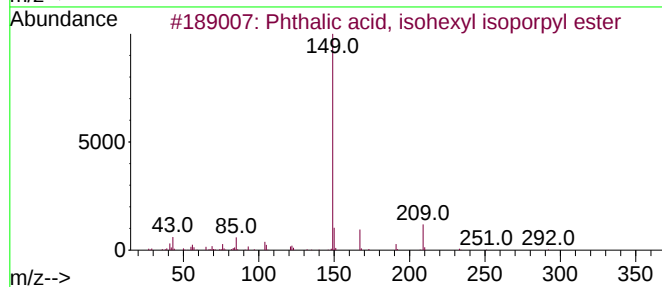
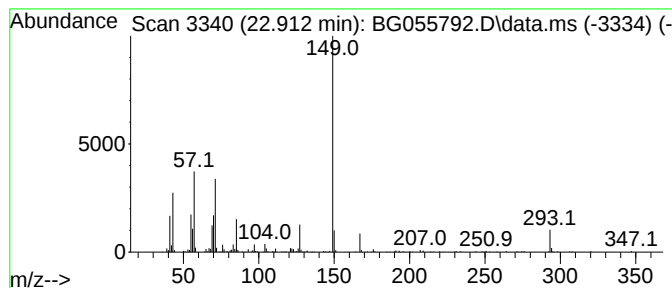
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Phthalic acid, isohexyl iso... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.913	15.92 ng	593782	Chrysene-d12	21.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, isohexyl isoporpy...	292	C17H24O4	1000315-00-4	64
2			Phthalic acid, 1-cyclopentylethy...	388	C24H36O4	1010309-73-8	59
3			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	59
4			Phthalic acid, 5-methylhex-2-yl ...	502	C32H54O4	1000371-07-7	58
5			Diamyl phthalate	306	C18H26O4	000131-18-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
Data File : BG055792.D
Acq On : 25 Nov 2022 22:48
Operator : CG/JU
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SOIL-PILE-10

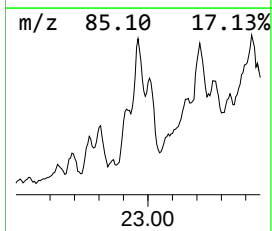
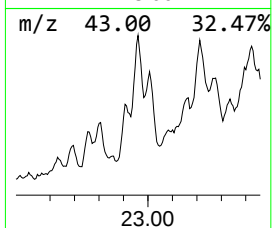
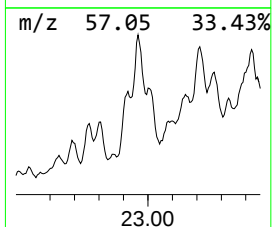
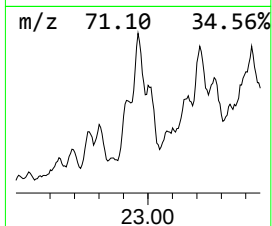
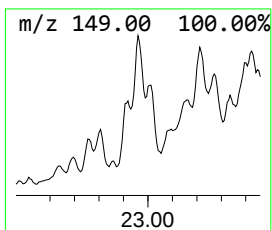
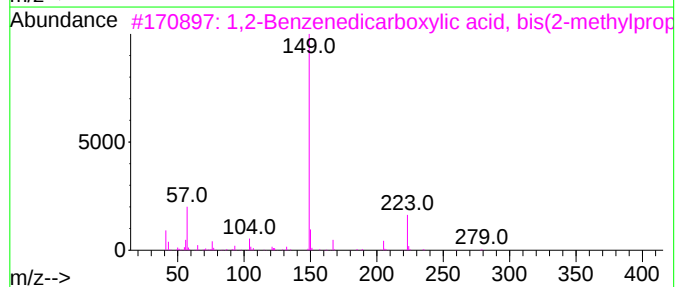
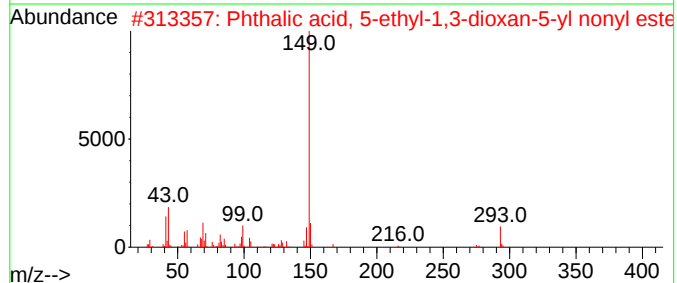
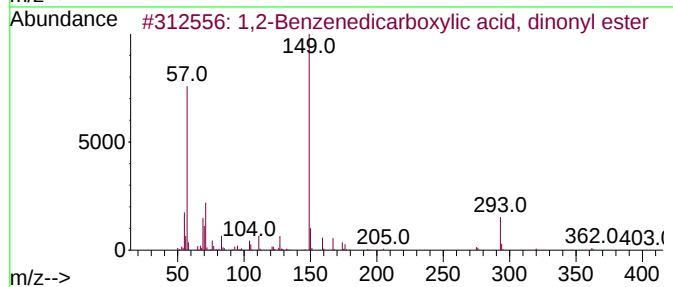
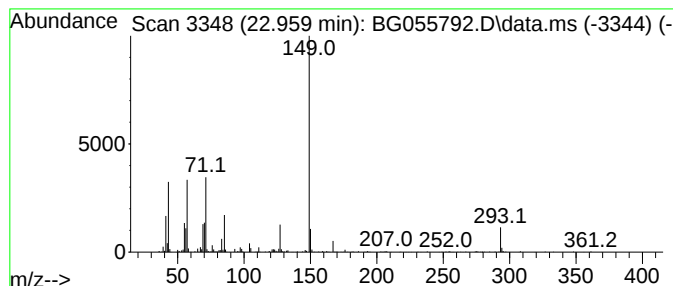
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1,2-Benzenedicarboxylic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.959	18.14 ng	676810	Chrysene-d12	21.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	72
2			Phthalic acid, 5-ethyl-1,3-dioxa...	420	C24H36O6	1000415-48-5	58
3			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	53
4			Phthalic acid, isobutyl nonyl ester	348	C21H32O4	1000309-04-4	50
5			Phthalic acid, pentyl 2-pentyl e...	306	C18H26O4	1000315-47-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
Data File : BG055792.D
Acq On : 25 Nov 2022 22:48
Operator : CG/JU
Sample : N5641-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
SOIL-PILE-10

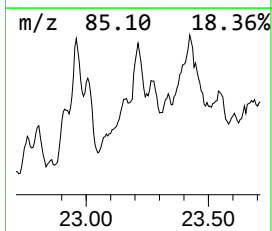
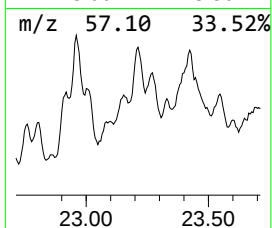
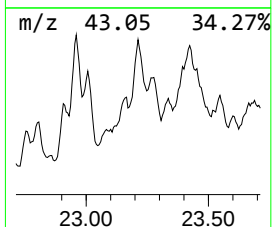
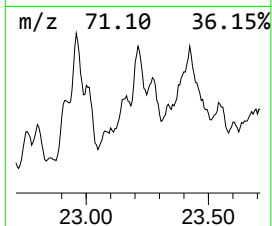
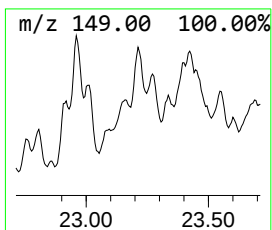
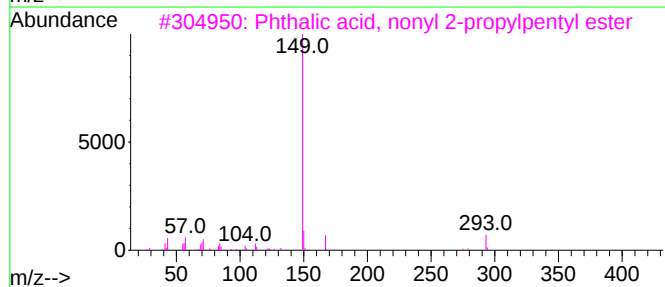
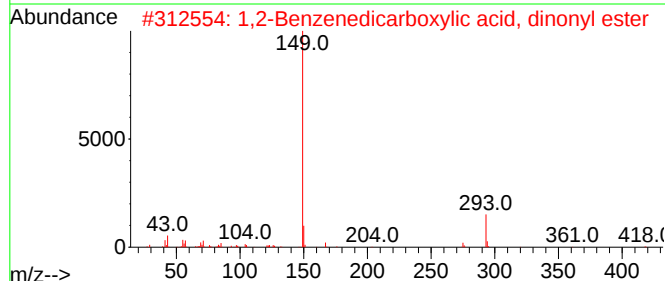
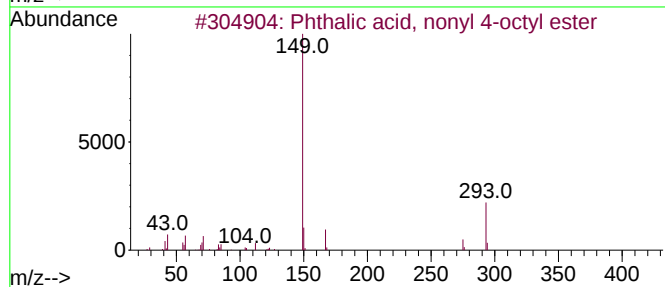
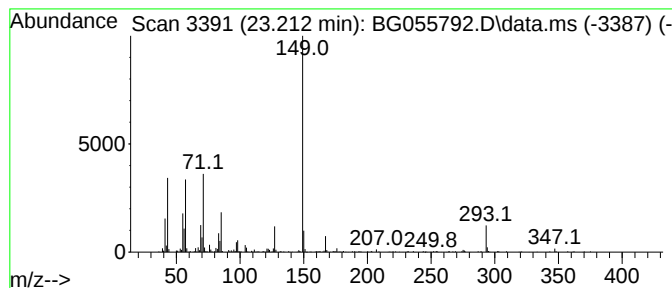
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Phthalic acid, nonyl 4-octyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.212	12.33 ng	459904	Chrysene-d12	21.890

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, nonyl 4-octyl ester	404	C25H40O4	1000314-84-2	72
2			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	59
3			Phthalic acid, nonyl 2-propylpen...	404	C25H40O4	1000377-92-5	59
4			Phthalic acid, hept-3-yl nonyl e...	390	C24H38O4	1000356-99-3	59
5			Phthalic acid, isohexyl 3-methyl...	320	C19H28O4	1000356-80-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112522\
Data File : BG055792.D
Acq On : 25 Nov 2022 22:48
Operator : CG/JU
Sample : N5641-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Ethanol, 2-(2-b...	10.803	9.0	ng	179983	2	11.062	400649	20.0
Tris(2-chloropr...	17.325	3.0	ng	94912	4	17.601	642812	20.0
n-Hexadecanoic ...	18.453	3.5	ng	113720	4	17.601	642812	20.0
Hexanoic acid, ...	21.297	2.1	ng	78307	5	21.890	746085	20.0
unknown21.502	21.502	2.0	ng	76213	5	21.890	746085	20.0
Phthalic acid, ...	22.695	4.2	ng	156729	5	21.890	746085	20.0
Phthalic acid, ...	22.760	7.8	ng	291096	5	21.890	746085	20.0
Phthalic acid, ...	22.913	15.9	ng	593782	5	21.890	746085	20.0
1,2-Benzenedica...	22.959	18.1	ng	676810	5	21.890	746085	20.0
Phthalic acid, ...	23.212	12.3	ng	459904	5	21.890	746085	20.0